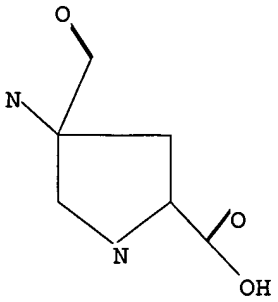
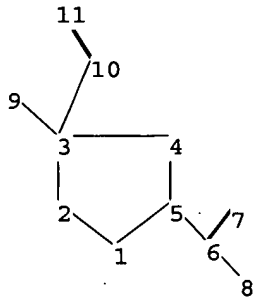
=>
Uploading C:\Program Files\Stnexp\Queries\10613961.str



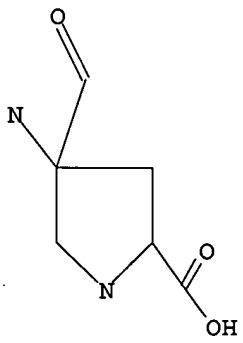


chain nodes :
6 7 8 9 10 11
ring nodes :
1 2 3 4 5
chain bonds :
3-9 3-10 5-6 6-7 6-8 10-11
ring bonds :
1-2 1-5 2-3 3-4 4-5
exact/norm bonds :
1-2 1-5 3-9 10-11
exact bonds :
2-3 3-4 3-10 4-5 5-6
normalized bonds :
6-7 6-8
isolated ring systems :
containing 1 :

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:CLASS 7:CLASS 8:CLASS 9:CLASS 10:CLASS
11:CLASS

L1 STRUCTURE UPLOADED

=> d l1
L1 HAS NO ANSWERS
L1 STR



Structure attributes must be viewed using STN Express query preparation.

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=> s l1
SAMPLE SEARCH INITIATED 15:24:00 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED -       14 TO ITERATE

100.0% PROCESSED       14 ITERATIONS                                 8 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS:  ONLINE  **COMPLETE**
                        BATCH   **COMPLETE**
PROJECTED ITERATIONS:           56 TO      504
PROJECTED ANSWERS:               8 TO      329

L2             8 SEA SSS SAM L1

=> s l1 ful
FULL SEARCH INITIATED 15:24:05 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED -      351 TO ITERATE

100.0% PROCESSED      351 ITERATIONS                             172 ANSWERS
SEARCH TIME: 00.00.01

L3           172 SEA SSS FUL L1

=> file caplus
COST IN U.S. DOLLARS                              SINCE FILE      TOTAL
                                                     ENTRY    SESSION
FULL ESTIMATED COST                                 166.94     167.15

FILE 'CAPLUS' ENTERED AT 15:24:07 ON 08 FEB 2006
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2006 AMERICAN CHEMICAL SOCIETY (ACS)
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FILE COVERS 1907 - 8 Feb 2006  VOL 144 ISS 7
FILE LAST UPDATED: 7 Feb 2006  (20060207/ED)

Effective October 17, 2005, revised CAS Information Use Policies apply.
They are available for your review at:

http://www.cas.org/infopolicy.html

=> s l3
L4             61 L3
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=> s l4 and py<2003
      22790841 PY<2003
L5           44 L4 AND PY<2003

=> s l5 and carbamate
          29354 CARBAMATE
          11121 CARBAMATES
          34008 CARBAMATE
                (CARBAMATE OR CARBAMATES)
L6            1 L5 AND CARBAMATE

=> d abs bib hitstr
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L6 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2006 ACS on STN

AB The four isomers of 4-aminopyrrolidine-2,4-dicarboxylate (APDC) were prepared and evaluated for their effects at glutamate receptors in vitro. (2R,4R)-APDC (2a), an aza analog of the nonselective mGluR agonist (1S,3R)-1-aminocyclopentane-1,3-dicarboxylate ((1S,3R)-ACPD, 1), was found to possess relatively high affinity for metabotropic glutamate receptors (mGluRs) (ACPD-sensitive [3H]glutamate binding IC50 = 6.49±1.21 μM) with no effects on radioligand binding to NMDA, AMPA, or kainate receptors up to 100 μM. None of the other APDC isomers showed significant mGluR binding affinity, indicating that this interaction is highly stereospecific. Both 1 and 2a were effective in decreasing forskolin-stimulated cAMP formation in the adult rat cerebral cortex (EC50 = 8.17±2.21 μM for 1; EC50 = 14.51±5.54 μM for 2a); however, while 1 was also effective in stimulating basal tritiated inositol monophosphate production in the neonatal rat cerebral cortex (EC50 = 27.7±5.2 μM), 2a (up to 100 μM) was ineffective in stimulating phosphoinositide hydrolysis in this tissue preparation, further supporting our previous observations that 2a is a highly selective agonist for mGluRs neg. coupled to adenylate cyclase. Microelectrophoretic application of either 1 or 2a to intact rat spinal neurons produced an augmentation of AMPA-induced excitation (95±10% increase for 1, 52±6% increase for 2a). Intracerebral injection of 1 (400 nmol) produced characteristic limbic seizures in mice which are not mimicked by 2a (200-1600 nmol, ic). However, the limbic seizures induced by 1 were blocked by systemically administered 2a in a dose-dependent manner (EC50 = 271 mg/kg, i.p.). It is concluded that (2R,4R)-APDC (2a) is a highly selective, systemically-active agonist of mGluRs neg. coupled to adenylate cyclase and that selective activation of these receptors in vivo can result in anticonvulsant effects.

AN 1996:383040 CAPLUS

DN 125:104243

TI Synthesis of the Four Isomers of 4-Aminopyrrolidine-2,4-dicarboxylate: Identification of a Potent, Highly Selective, and Systemically-Active Agonist for Metabotropic Glutamate Receptors Negatively Coupled to Adenylate Cyclase

AU Monn, James A.; Valli, Matthew J.; Johnson, Bryan G.; Salhoff, Craig R.; Wright, Rebecca A.; Howe, Trevor; Bond, Ann; Lodge, David; Spangle, Larry A.; et al.

CS Core Technology Division, Eli Lilly and Company, Indianapolis, IN, USA

SO Journal of Medicinal Chemistry (**1996**), 39(15), 2990-3000
CODEN: JMCMAR; ISSN: 0022-2623

PB American Chemical Society

DT Journal

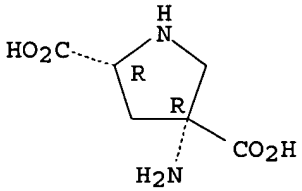
LA English

OS CASREACT 125:104243

IT **169209-63-6P**, (2R,4R)-4-Aminopyrrolidine-2,4-dicarboxylate
171336-79-1P 171336-80-4P 171336-81-5P
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); PRP (Properties); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(synthesis of four isomers of 4-aminopyrrolidine-2,4-dicarboxylate as agonists for metabotropic glutamate receptors neg. coupled to adenylate cyclase)

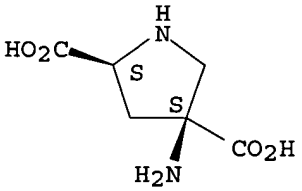
RN 169209-63-6 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-, (2R,4R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (+).



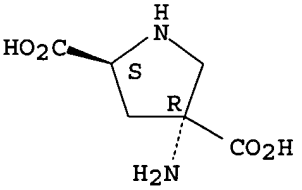
RN 171336-79-1 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-, (2S-cis)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (+).



RN 171336-80-4 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-, (2S-trans)- (9CI) (CA INDEX NAME)

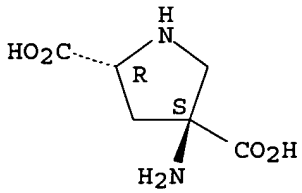
Absolute stereochemistry. Rotation (-).



RN 171336-81-5 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-, (2R-trans)- (9CI) (CA INDEX

NAME)

Absolute stereochemistry. Rotation (+).



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       1516378 GROUP
        988533 GROUPS
       2120181 GROUP
                 (GROUP OR GROUPS)
         15139 PROTECTING GROUP
                 (PROTECTING(W)GROUP)
L7           1 L5 AND PROTECTING GROUP

=> d l7 not l6
L6 IS NOT VALID HERE
For an explanation, enter "HELP DISPLAY".

=> d abs l7

L7   ANSWER 1 OF 1  CAPLUS  COPYRIGHT 2006 ACS on STN
GI
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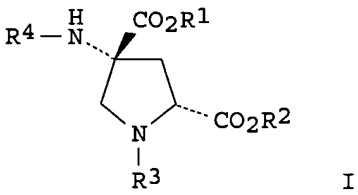


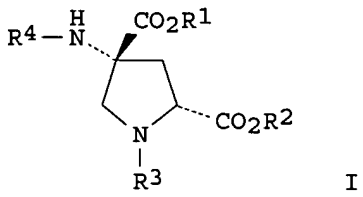
I

AB The present invention provides pyrrolidinyl dicarboxylic acid derivs. I wherein: R1 and R2 are each individually H or a carboxy **protecting group**; R4 is H or an amino **protecting group**; R3 = e.g., C1-16 alkyl, C3-8 cycloalkyl; C3-8 cycloalkenyl, aryl, that affect certain excitatory amino acid receptors (no data), and are useful in the treatment of neurol. disorders and psychiatric disorders. This invention further provides novel pyrrolidinyl di-carboxylic acid derivs. and pharmaceutical formulations employing these novel compds. Thus, cis-4-hydroxy-D-proline was esterified and N-benzylated to provide (2R,4R) Et 1-benzyl-4-hydroxypyrrolidine-2-carboxylate; this was oxidized to the 4-oxo derivative which was treated with KCN/ammonium carbonate to afford (2R,4R) di-Et 1-benzyl-4-aminopyrrolidine-2,4-dicarboxylate; the latter

was N-protected and debenzylated to afford (2R,4R) di-Et 4-(BOC-amino)pyrrolidine-2,4-dicarboxylate (II) as the scaffold intermediate. Reductive alkylation of II with pentanal afforded (2R,4R) di-Et 4-(BOC-amino)-1-pentylpyrrolidine-2,4-dicarboxylate which was deprotected and hydrolyzed to (2R,4R) 4-amino-1-pentylpyrrolidine-2,4-dicarboxylic acid (I; R1 = R2 = R4 = H, R3 = pentyl).

=> d abs fbib hitstr

L7 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2006 ACS on STN
GI



I

AB The present invention provides pyrrolidinyl dicarboxylic acid derivs. I wherein: R1 and R2 are each individually H or a carboxy **protecting group**; R4 is H or an amino **protecting group**; R3 = e.g., C1-16 alkyl, C3-8 cycloalkyl; C3-8 cycloalkenyl, aryl, that affect certain excitatory amino acid receptors (no data), and are useful in the treatment of neurol. disorders and psychiatric disorders. This invention further provides novel pyrrolidinyl di-carboxylic acid derivs. and pharmaceutical formulations employing these novel compds. Thus, cis-4-hydroxy-D-proline was esterified and N-benzylated to provide (2R,4R) Et 1-benzyl-4-hydroxypyrrolidine-2-carboxylate; this was oxidized to the 4-oxo derivative which was treated with KCN/ammonium carbonate to afford (2R,4R) di-Et 1-benzyl-4-aminopyrrolidine-2,4-dicarboxylate; the latter was N-protected and debenzylated to afford (2R,4R) di-Et 4-(BOC-amino)pyrrolidine-2,4-dicarboxylate (II) as the scaffold intermediate. Reductive alkylation of II with pentanal afforded (2R,4R) di-Et 4-(BOC-amino)-1-pentylpyrrolidine-2,4-dicarboxylate which was deprotected and hydrolyzed to (2R,4R) 4-amino-1-pentylpyrrolidine-2,4-dicarboxylic acid (I; R1 = R2 = R4 = H, R3 = pentyl).
AN 1996:410401 CAPLUS
DN 125:86486
TI (2R,4R)-4-Aminopyrrolidine-2,4-dicarboxylic acid derivatives as metabotropic glutamate receptor antagonists
IN Monn, James Allen; Tizzano, Joseph Patrick; Valli, Matthew J.
PA Eli Lilly and Co., USA
SO PCT Int. Appl., 97 pp.
CODEN: PIXXD2
DT Patent
LA English
FAN.CNT 1

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI WO 9605828	A1	19960229	WO 1995-US10320	19950814 <--

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MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, TJ, TM
RW: KE, MW, SD, SZ, UG, AT, BE, CH, DE, DK, ES, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, ML, MR, NE, SN, TD, TG

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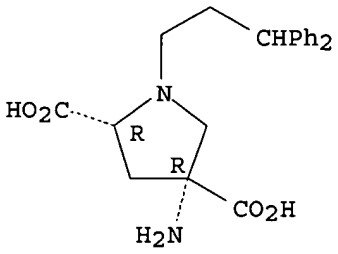
OS MARPAT 125:86486

IT **178415-98-0P**
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); RCT (Reactant); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent); USES (Uses)
((2R,4R)-4-aminopyrrolidine-2,4-dicarboxylic acid derivs. as metabotropic glutamate receptor antagonists)

RN 178415-98-0 CAPLUS

CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-(3,3-diphenylpropyl)-, (2R,4R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



IT **171336-76-8P 178415-42-4P 178415-44-6P**
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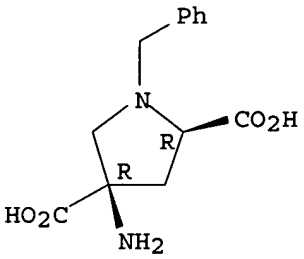
178416-01-8P 178416-03-0P 178416-05-2P
178416-07-4P 178416-09-6P 178416-11-0P
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
((2R,4R)-4-aminopyrrolidine-2,4-dicarboxylic acid derivs. as

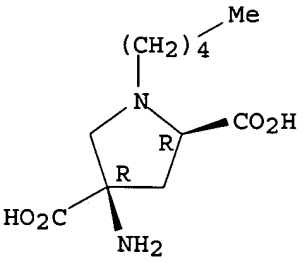
metabotropic glutamate receptor antagonists)
RN 171336-76-8 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-(phenylmethyl)-, (2R,4R)-(9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (+).



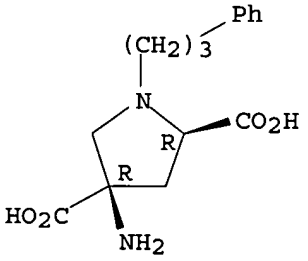
RN 178415-42-4 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-pentyl-, (2R-cis)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



RN 178415-44-6 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-(3-phenylpropyl)-, (2R,4R)-(9CI) (CA INDEX NAME)

Absolute stereochemistry.



RN 178415-46-8 CAPLUS
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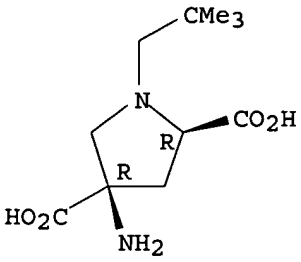
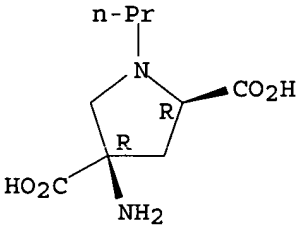
Absolute stereochemistry.



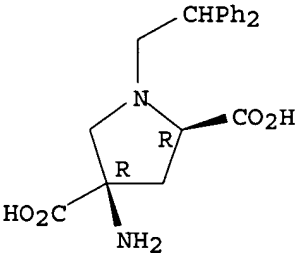
RN 178415-48-0 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-propyl-, (2R-cis)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



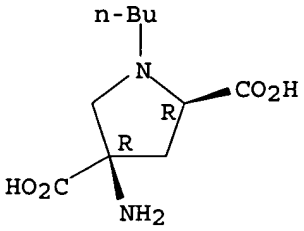
RN 178415-50-4 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-(2,2-diphenylethyl)-, (2R,4R)-(9CI) (CA INDEX NAME)

Absolute stereochemistry.



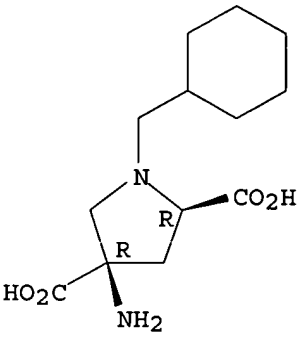
RN 178415-52-6 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-butyl-, (2R-cis)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



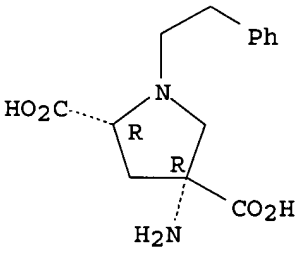
RN 178415-54-8 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-(cyclohexylmethyl)-, (2R,4R)-(9CI) (CA INDEX NAME)

Absolute stereochemistry.



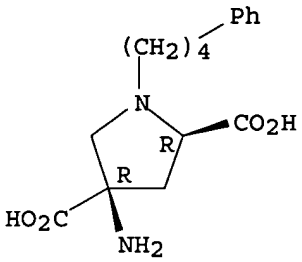
RN 178415-56-0 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-(2-phenylethyl)-, (2R,4R)-(9CI) (CA INDEX NAME)

Absolute stereochemistry.



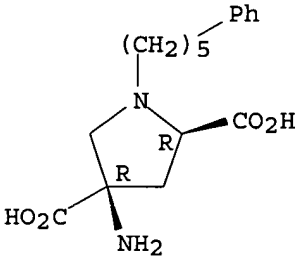
RN 178415-58-2 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-(4-phenylbutyl)-, (2R-cis)-(9CI) (CA INDEX NAME)

Absolute stereochemistry.



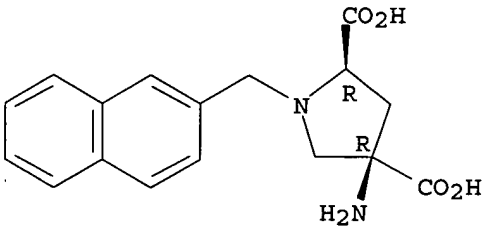
RN 178415-60-6 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-(5-phenylpentyl)-, (2R-cis)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



RN 178415-62-8 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-(2-naphthalenylmethyl)-, (2R,4R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



RN 178415-64-0 CAPLUS
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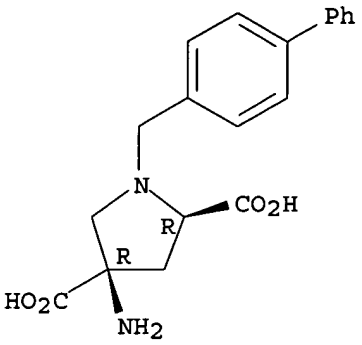
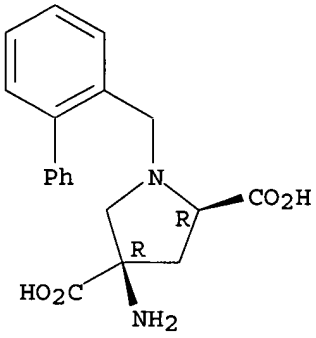
Absolute stereochemistry.



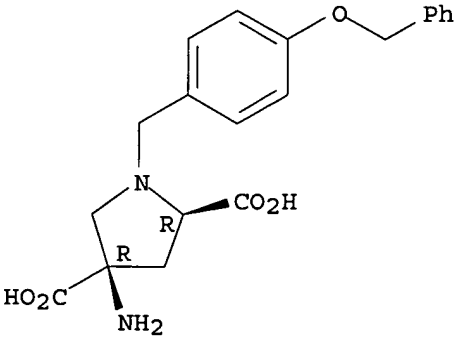
RN 178415-66-2 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-([1,1'-biphenyl]-2-ylmethyl)-, (2R,4R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



RN 178415-68-4 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-[[4-(phenylmethoxy)phenyl]methyl]-, (2R-cis)- (9CI) (CA INDEX NAME)

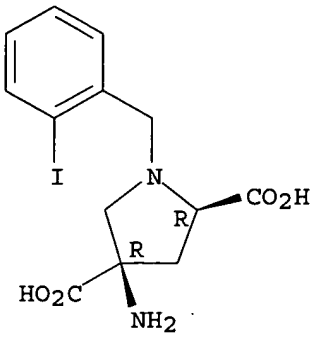
Absolute stereochemistry.



RN 178415-70-8 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-[(2-iodophenyl)methyl]-,

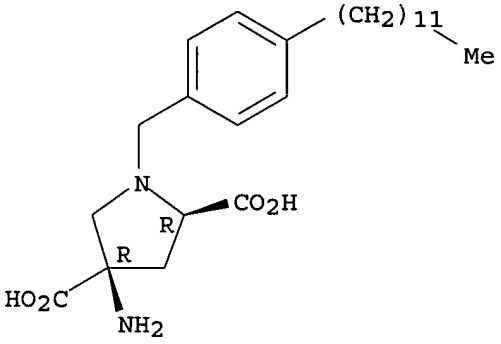
(2R-cis)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



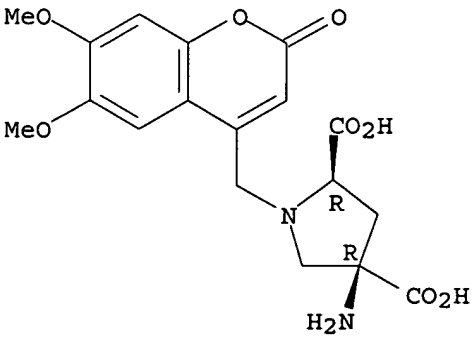
RN 178415-72-0 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-[(4-dodecylphenyl)methyl]-, (2R-cis)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



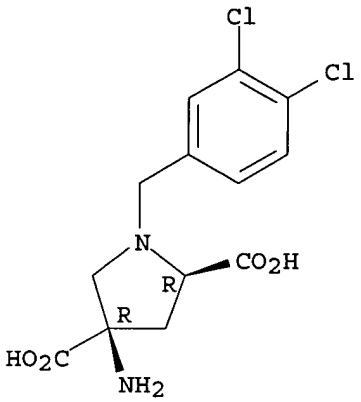
RN 178415-74-2 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-[(6,7-dimethoxy-2-oxo-2H-1-benzopyran-4-yl)methyl]-, (2R-cis)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



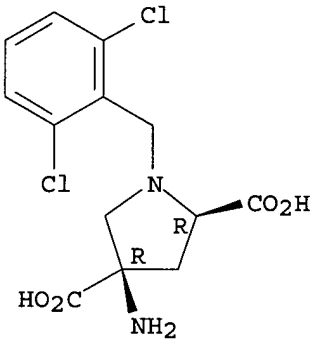
RN 178415-76-4 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-[(3,4-dichlorophenyl)methyl]-, (2R,4R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



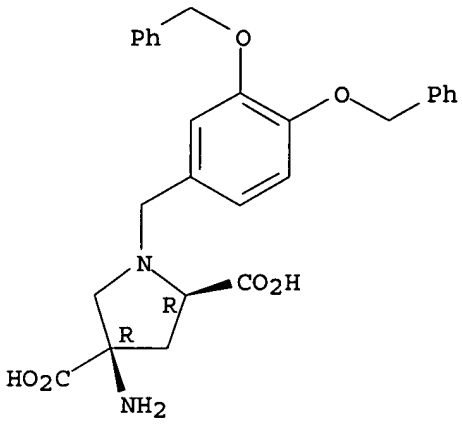
RN 178415-78-6 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-[(2,6-dichlorophenyl)methyl]-, (2R,4R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



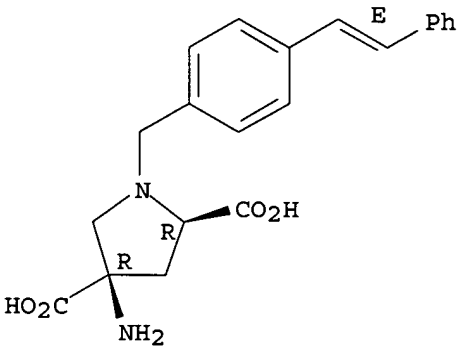
RN 178415-80-0 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-[[3,4-bis(phenylmethoxy)phenyl]methyl]-, (2R-cis)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



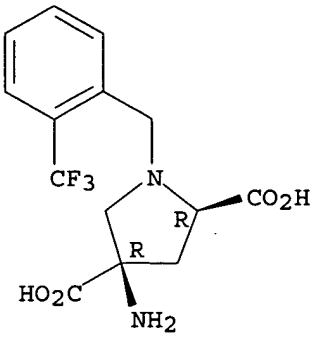
RN 178415-82-2 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-[[4-(2-phenylethenyl)phenyl]methyl]-, [2R-[1(E),2α,4α]]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.
Double bond geometry as shown.



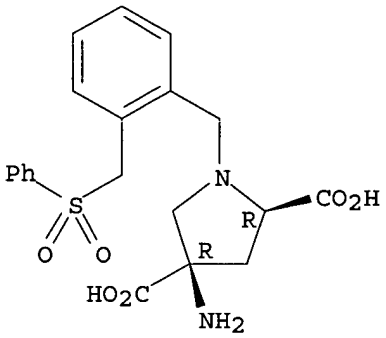
RN 178415-84-4 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-[[2-(trifluoromethyl)phenyl]methyl]-, (2R-cis)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



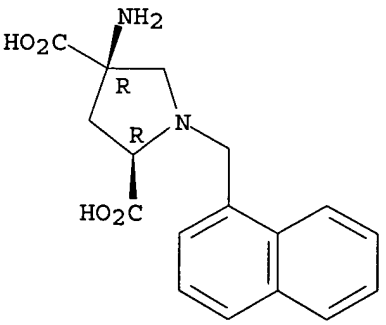
RN 178415-86-6 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-[[2-[(phenylsulfonyl)methyl]phenyl]methyl]-, (2R-cis)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



RN 178415-88-8 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-(1-naphthalenylmethyl)-, (2R,4R)- (9CI) (CA INDEX NAME)

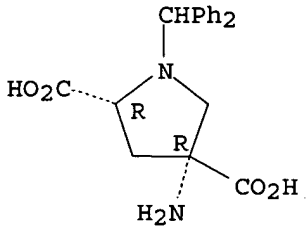
Absolute stereochemistry.



RN 178415-90-2 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-(diphenylmethyl)-, (2R,4R)-

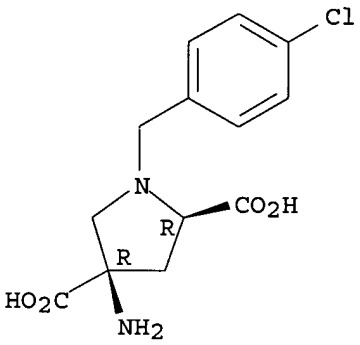
(9CI) (CA INDEX NAME)

Absolute stereochemistry.



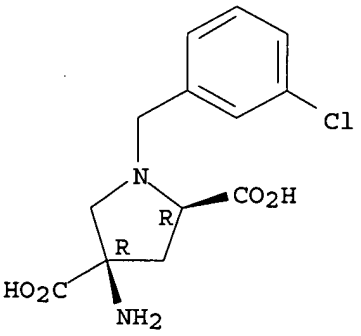
RN 178415-92-4 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-[(4-chlorophenyl)methyl]-, (2R,4R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



RN 178415-94-6 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-[(3-chlorophenyl)methyl]-, (2R,4R)- (9CI) (CA INDEX NAME)

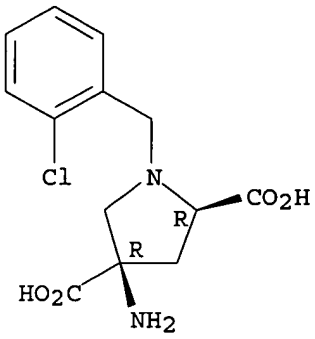
Absolute stereochemistry.



RN 178415-96-8 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-[(2-chlorophenyl)methyl]-,

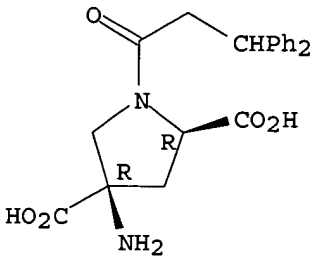
(2R,4R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



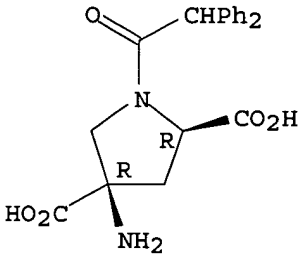
RN 178416-00-7 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-(1-oxo-3,3-diphenylpropyl)-, (2R-cis)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



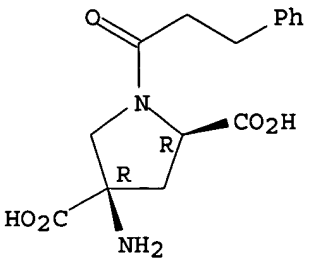
RN 178416-01-8 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-(diphenylacetyl)-, (2R-cis)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



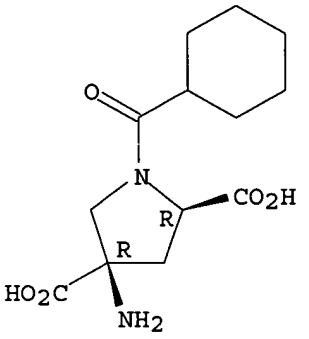
RN 178416-03-0 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-(1-oxo-3-phenylpropyl)-, (2R-cis)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



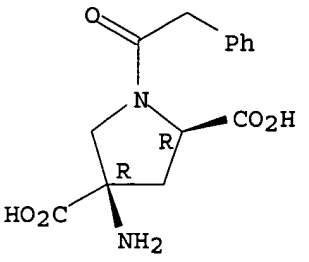
RN 178416-05-2 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-(cyclohexylcarbonyl)-, (2R-cis)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



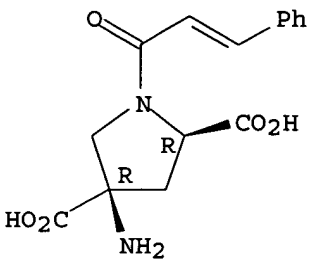
RN 178416-07-4 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-(phenylacetyl)-, (2R-cis)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



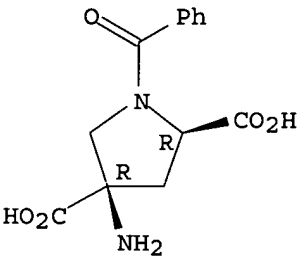
RN 178416-09-6 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-(1-oxo-3-phenyl-2-propenyl)-, (2R-cis)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.
Double bond geometry unknown.



RN 178416-11-0 CAPLUS
CN 2,4-Pyrrolidinedicarboxylic acid, 4-amino-1-benzoyl-, (2R-cis)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



=> file stnguide

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	23.76	190.91

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	-2.25	-2.25

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FILE CONTAINS CURRENT INFORMATION.
LAST RELOADED: Feb 3, 2006 (20060203/UP).

=> d his

(FILE 'HOME' ENTERED AT 15:23:30 ON 08 FEB 2006)

FILE 'REGISTRY' ENTERED AT 15:23:40 ON 08 FEB 2006
L1 STRUCTURE UPLOADED
L2 8 S L1
L3 172 S L1 FUL

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     FILE 'CAPLUS' ENTERED AT 15:24:07 ON 08 FEB 2006
L4            61 S L3
L5            44 S L4 AND PY<2003
L6             1 S L5 AND CARBAMATE
L7             1 S L5 AND PROTECTING GROUP

     FILE 'STNGUIDE' ENTERED AT 15:29:33 ON 08 FEB 2006

=> s l4 and ?fluorenyl?
COMBINATION OF STRUCTURE AND TEXT TERMS NOT VALID
The query entered contains both search terms created by
structure-building or screen commands and text search terms.  L#s
created via the STRUCTURE or SCREEN commands must be searched in the
structures files separately from text terms or profiles.  The L#
answer sets from structure searches can be used in crossover searches
and can be combined with text terms.

=> s l4 and fluorenyl
COMBINATION OF STRUCTURE AND TEXT TERMS NOT VALID
The query entered contains both search terms created by
structure-building or screen commands and text search terms.  L#s
created via the STRUCTURE or SCREEN commands must be searched in the
structures files separately from text terms or profiles.  The L#
answer sets from structure searches can be used in crossover searches
and can be combined with text terms.

=> s l5 and fluorenyl
COMBINATION OF STRUCTURE AND TEXT TERMS NOT VALID
The query entered contains both search terms created by
structure-building or screen commands and text search terms.  L#s
created via the STRUCTURE or SCREEN commands must be searched in the
structures files separately from text terms or profiles.  The L#
answer sets from structure searches can be used in crossover searches
and can be combined with text terms.

=> file caplus
COST IN U.S. DOLLARS                            SINCE FILE      TOTAL
                                                   ENTRY    SESSION
FULL ESTIMATED COST                                 0.78     191.69

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)      SINCE FILE      TOTAL
                                                   ENTRY    SESSION
CA SUBSCRIBER PRICE                                 0.00      -2.25

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FILE LAST UPDATED: 7 Feb 2006 (20060207/ED)

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=> s l5 and fluorenyl
          4084 FLUORENYL
             5 FLUORENYLS
          4085 FLUORENYL
                 (FLUORENYL OR FLUORENYLS)
L8           0 L5 AND FLUORENYL

=> s l5 and fluorenyl?
          9942 FLUORENYL?
L9           0 L5 AND FLUORENYL?
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